LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.273	26	31	36	rBV2	4629	5850	0.80%	0.088%
2	1.399	65	70	78	rBV	94783	98642	13.45%	1.478%
3	1.681	151	158	178	rVB	73873	97436	13.29%	1.460%
4	2.273	332	342	354	rBV	147144	223194	30.44%	3.344%
5	2.443	388	395	396	rBV2	1650	1473	0.20%	0.022%
6	2.566	430	433	436	rBV	246	179	0.02%	0.003%
7	2.588	437	440	443	rBV	265	212	0.03%	0.003%
8	2.842	516	519	522	rVB2	235	184	0.03%	0.003%
9	3.189	620	627	637	rVB4	1169	2144	0.29%	0.032%
10	3.382	685	687	691	rVB2	342	192	0.03%	0.003%
11	4.180	922	935	959	rBV2	76487	198916	27.13%	2.980%
12	4.495	1029	1033	1038	rVB2	360	364	0.05%	0.005%
13	4.652	1064	1082	1109	rBV	119123	281907	38.45%	4.224%
14	4.881	1150	1153	1155	rBV	338	251	0.03%	0.004%
15	5.344	1273	1297	1335	rVV2	240687	733143	100.00%	10.985%
16	5.501	1345	1346	1354	rVB2	379	333	0.05%	0.005%
17	5.781	1427	1433	1434	rBV2	1103	804	0.11%	0.012%
18	5.890	1453	1467	1503	rBV	235763	473528	64.59%	7.095%
19	6.170	1550	1554	1556	rBV3	419	292	0.04%	0.004%
20	6.334	1591	1605	1639	rVV	172315	350851	47.86%	5.257%
21	6.829	1749	1759	1774	rBV3	6025	11316	1.54%	0.170%
22	7.231	1872	1884	1902	rBV	92194	167306	22.82%	2.507%
23	7.395	1924	1935	1946	rBV4	6227	12047	1.64%	0.181%
24	7.485	1958	1963	1968	rBV2	154	203	0.03%	0.003%
25	7.569	1976	1989	2002	rVV	304847	534743	72.94%	8.012%
26	7.639	2003	2011	2029	rVB	42104	77314	10.55%	1.158%
27	7.752	2043	2046	2048	rBV2	251	186	0.03%	0.003%
28	7.852	2063	2077	2102	rBV	67005	115716	15.78%	1.734%
29	7.951	2102	2108	2114	rVB5	612	818	0.11%	0.012%
30	8.028	2128	2132	2134	rBV	299	258	0.04%	0.004%
31	8.073	2142	2146	2147	rBV2	252	178	0.02%	0.003%
32	8.086	2147	2150	2155	rVB	412	276	0.04%	0.004%
33	8.183	2168	2180	2188	rBV2	4367	9780	1.33%	0.147%
34	8.311	2210	2220	2251	rVB	258341	452789	61.76%	6.784%

LSC Area Percent Report

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Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Title : VOC Analysis

35	8.572	2297	2301	2305	rBV2	284	285	0.04%	0.004%
36	8.649	2321	2325	2329	rBV2	230	194	0.03%	0.003%
37	8.733	2345	2351	2354	rBV	256	322	0.04%	0.005%
38	8.839	2381	2384	2389	rBV2	201	192	0.03%	0.003%
39	9.090	2449	2462	2488	rBV	350971	580960	79.24%	8.705%
40	9.254	2502	2513	2527	rBV2	9551	16082	2.19%	0.241%
41	9.379	2542	2552	2564	rBV	15982	26275	3.58%	0.394%
42	9.446	2569	2573	2582	rVV4	1121	1629	0.22%	0.024%
43	9.524	2591	2597	2600	rBV2	260	307	0.04%	0.005%
44	9.662	2635	2640	2642	rBV	236	218	0.03%	0.003%
45	9.749	2663	2667	2668	rBV2	256	183	0.02%	0.003%
46	9.781	2668	2677	2689	rVV	14310	22407	3.06%	0.336%
47	9.900	2707	2714	2719	rBV3	384	578	0.08%	0.009%
48	10.134	2784	2787	2791	rVB2	423	357	0.05%	0.005%
49	10.170	2792	2798	2803	rVV3	1232	1621	0.22%	0.024%
50	10.318	2840	2844	2852	rVB3	1106	975	0.13%	0.015%
51	10.434	2867	2880	2901	rVB	228207	365083	49.80%	5.470%
52	10.524	2901	2908	2913	rBV2	278	340	0.05%	0.005%
53	10.614	2920	2936	2949	rBV5	8134	19041	2.60%	0.285%
54	10.681	2949	2957	2959	rBV4	2012	2615	0.36%	0.039%
55	10.781	2981	2988	2992	rVB2	762	973	0.13%	0.015%
56	10.835	3002	3005	3009	rVB	265	243	0.03%	0.004%
57	10.913	3026	3029	3032	rBV	290	242	0.03%	0.004%
58	10.974	3039	3048	3053	rBV3	3928	5379	0.73%	0.081%
59	11.151	3094	3103	3115	rVB2	11732	17648	2.41%	0.264%
60	11.215	3115	3123	3129	rBV5	4025	5302	0.72%	0.079%
61	11.282	3138	3144	3150	rVB6	1176	1638	0.22%	0.025%
62	11.340	3155	3162	3163	rBV3	1072	1103	0.15%	0.017%
63	11.485	3192	3207	3225	rBV	367722	567924	77.46%	8.509%
64	11.572	3225	3234	3239	rBV5	3392	5483	0.75%	0.082%
65	11.633	3248	3253	3260	rVB7	3277	4313	0.59%	0.065%
66	11.765	3281	3294	3304	rBV2	24114	43343	5.91%	0.649%
67	11.861	3307	3324	3349	rVB	322422	517663	70.61%	7.756%
68	11.983	3351	3362	3369	rBV	24580	38102	5.20%	0.571%
69	12.122	3398	3405	3414	rBV3	2900	3787	0.52%	0.057%
70	12.212	3431	3433	3434	rBV	459	200	0.03%	0.003%
71	12.289	3450	3457	3463	rBV3	5560	6826	0.93%	0.102%

LSC Area Percent Report

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Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Title : VOC Analysis

72	12.356	3472	3478	3481	rBV2	4121	5024	0.69%	0.075%
73	12.388	3482	3488	3495	rVV5	12867	20598	2.81%	0.309%
74	12.437	3497	3503	3509	rVV6	6092	9815	1.34%	0.147%
75	12.478	3510	3516	3524	rVV5	8038	14079	1.92%	0.211%
76	12.523	3525	3530	3542	rVV5	6642	15185	2.07%	0.228%
77	12.594	3544	3552	3562	rVB4	17629	29422	4.01%	0.441%
78	12.697	3573	3584	3589	rBV4	32962	65356	8.91%	0.979%
79	12.742	3593	3598	3619	rVB	88294	154382	21.06%	2.313%
80	12.893	3641	3645	3652	rBV3	1166	1171	0.16%	0.018%
81	12.925	3652	3655	3656	rBV2	1235	625	0.09%	0.009%
82	12.964	3662	3667	3677	rVB5	5955	8353	1.14%	0.125%
83	13.093	3701	3707	3709	rBV4	1494	1621	0.22%	0.024%
84	13.128	3710	3718	3728	rVV6	13547	24435	3.33%	0.366%
85	13.186	3729	3736	3749	rVB2	11531	17707	2.42%	0.265%
86	13.292	3758	3769	3781	rVB4	21606	37130	5.06%	0.556%
87	13.392	3788	3800	3812	rVB10	2166	4447	0.61%	0.067%
88	13.456	3812	3820	3825	rBV7	5780	8742	1.19%	0.131%
89	13.620	3866	3871	3880	rBV7	2423	3882	0.53%	0.058%
90	13.675	3882	3888	3896	rVB8	2715	4382	0.60%	0.066%
91	13.742	3898	3909	3916	rBV	33460	61727	8.42%	0.925%
92	13.903	3951	3959	3969	rVB	27528	42376	5.78%	0.635%
93	14.006	3984	3991	4001	rVB6	6935	10268	1.40%	0.154%
94	14.054	4001	4006	4014	rVB7	2550	3573	0.49%	0.054%
95	14.218	4048	4057	4066	rVB7	3405	6550	0.89%	0.098%
96	14.253	4066	4068	4069	rBV	483	250	0.03%	0.004%
97	14.392	4104	4111	4114	rBV7	1373	1834	0.25%	0.027%
98	14.523	4150	4152	4154	rBV	427	236	0.03%	0.004%
99	14.671	4196	4198	4200	rBV2	473	310	0.04%	0.005%
100	15.067	4312	4321	4332	rBV	5266	7880	1.07%	0.118%

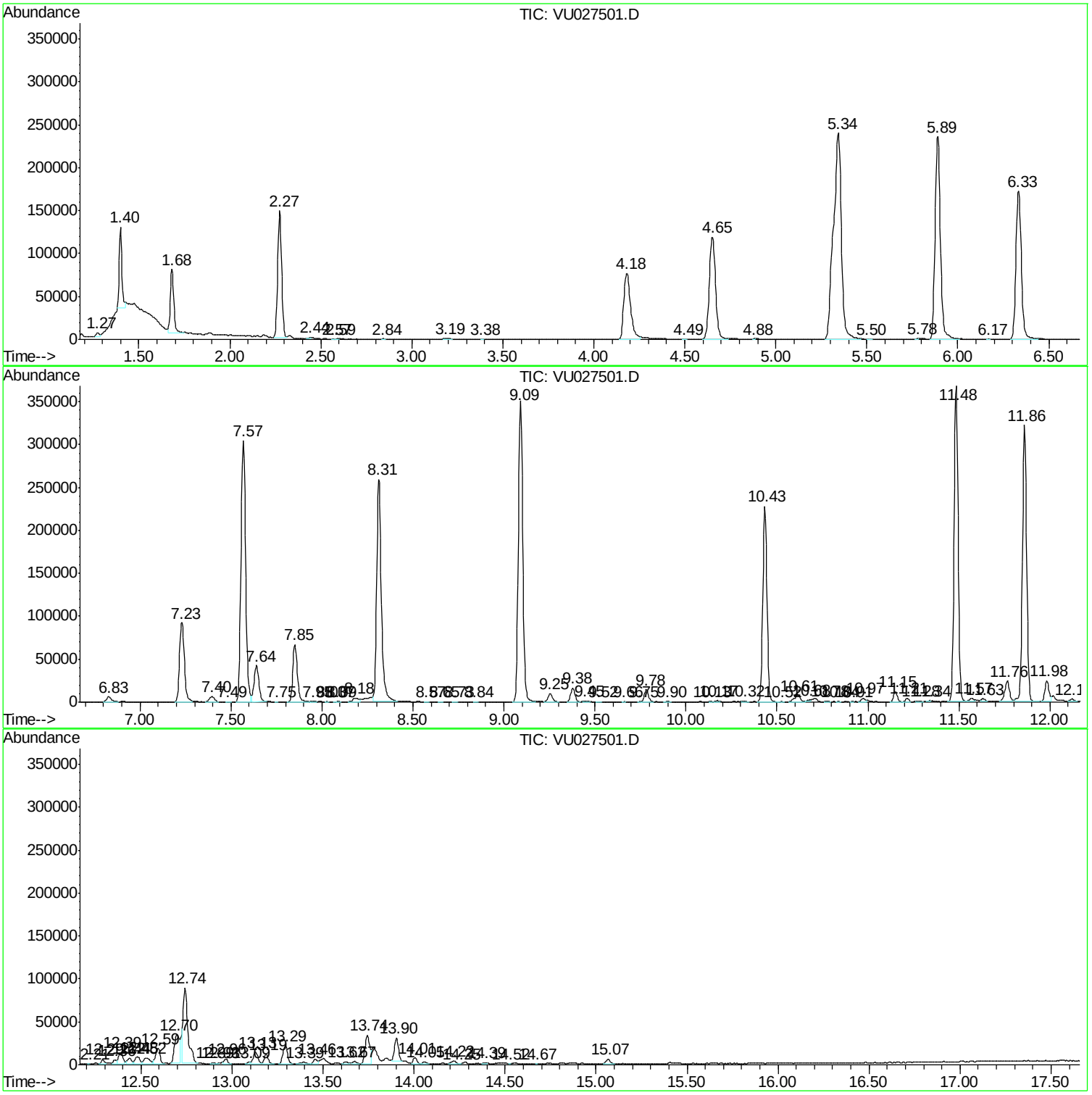
Sum of corrected areas: 6674018

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100818\
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Quant Title : VOC Analysis

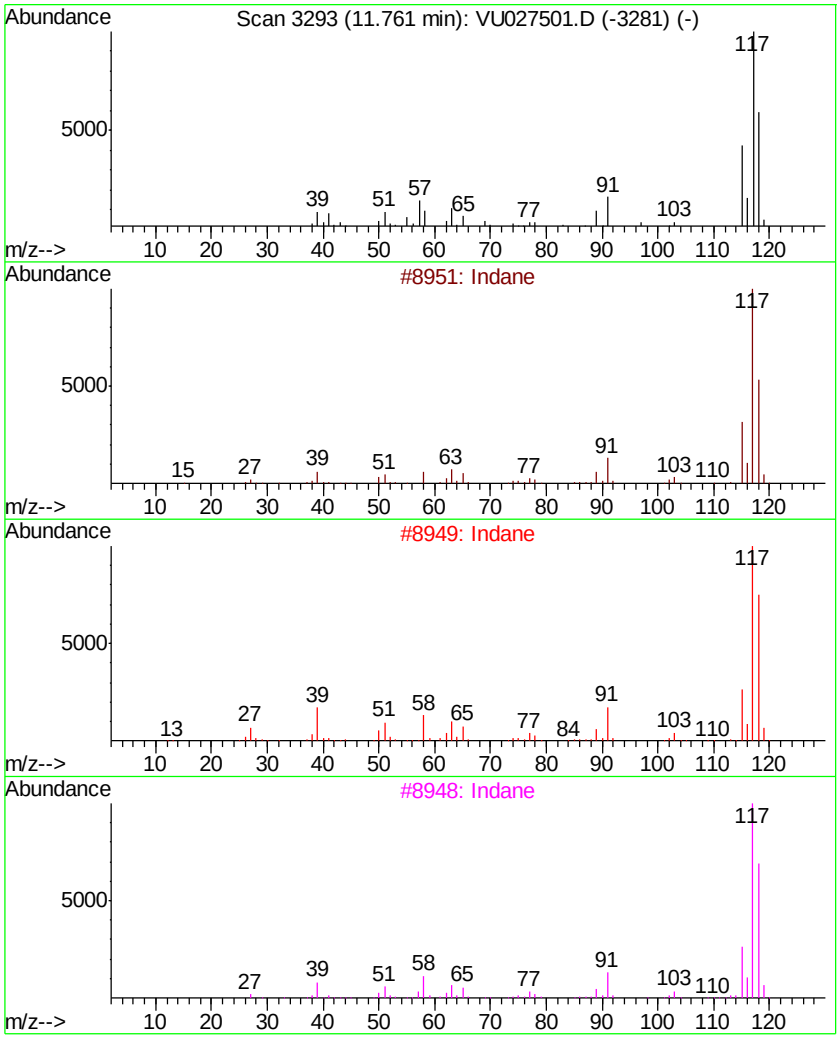
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

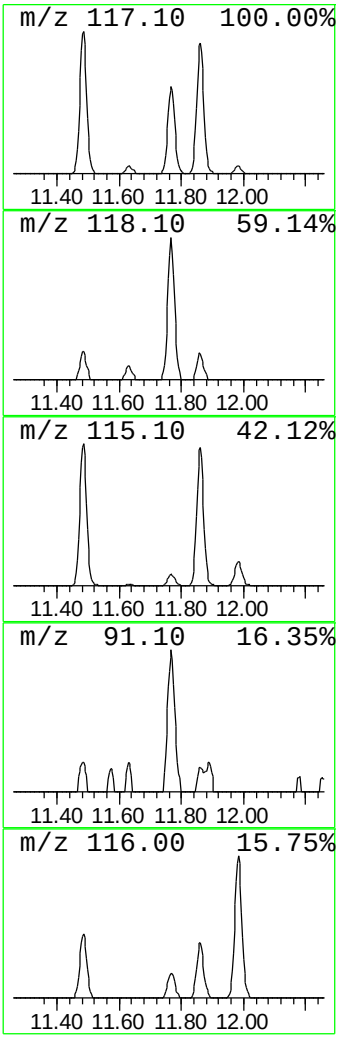
****

Peak Number 2 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	3.82 ug/L	43343	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Indane	118	C9H10	000496-11-7	94
3	Indane	118	C9H10	000496-11-7	70
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	62





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

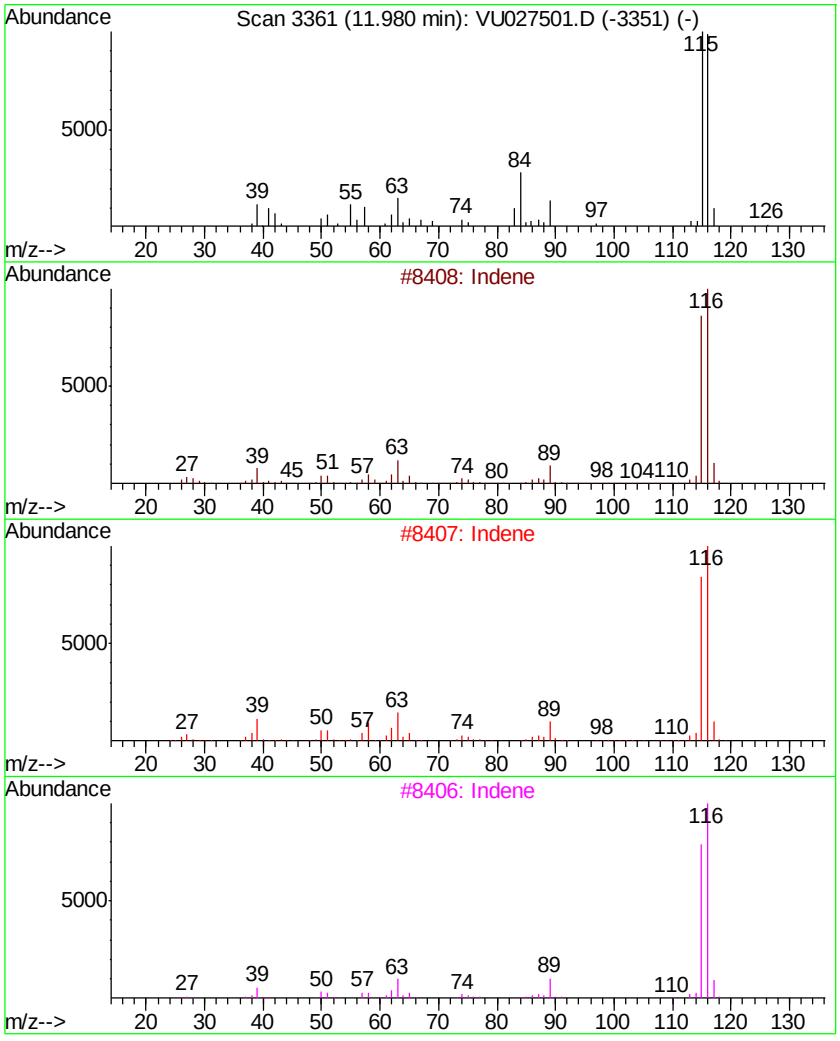
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

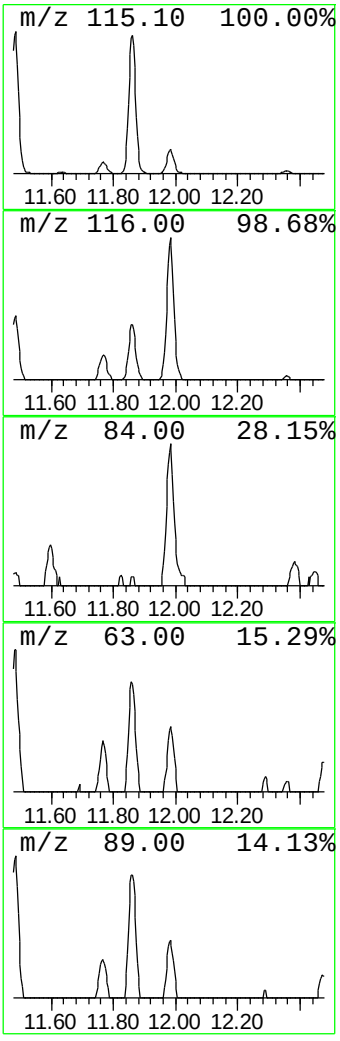
****

Peak Number 3 Indene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	3.35 ug/L	38102	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	95
2	Indene	116	C9H8	000095-13-6	93
3	Indene	116	C9H8	000095-13-6	91
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
5	3-Methylphenylacetylene	116	C9H8	000766-82-5	90





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

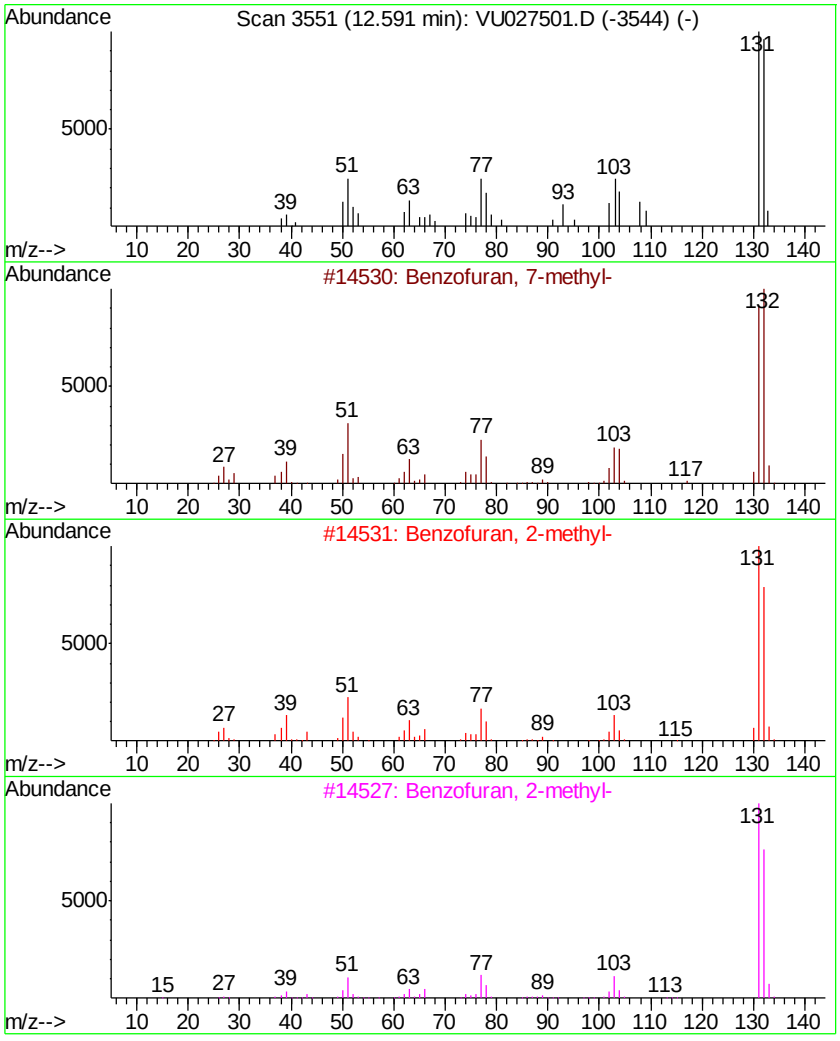
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

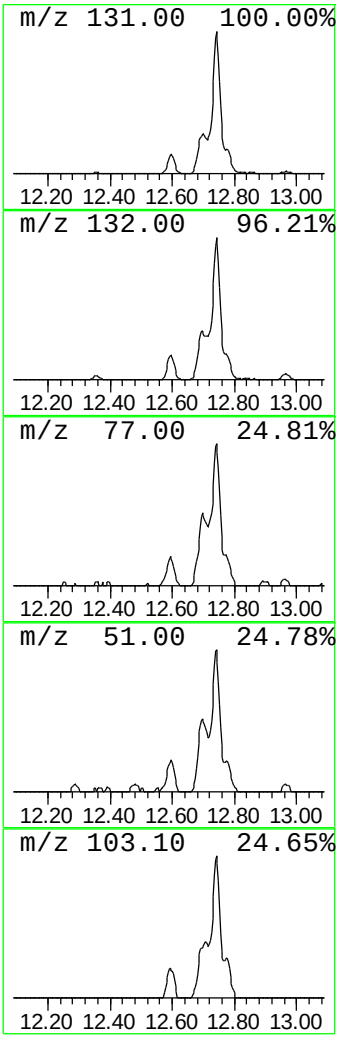
****

Peak Number 4 Benzofuran, 7-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	2.59 ug/L	29422	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	95
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
4	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

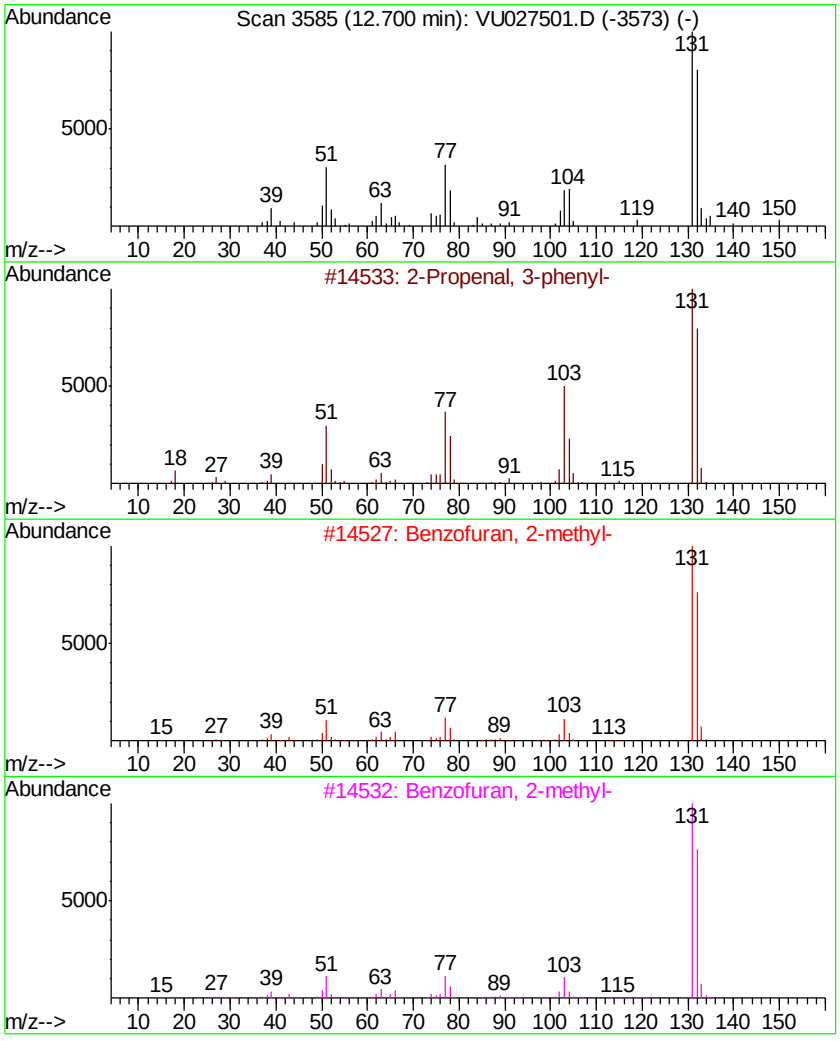
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

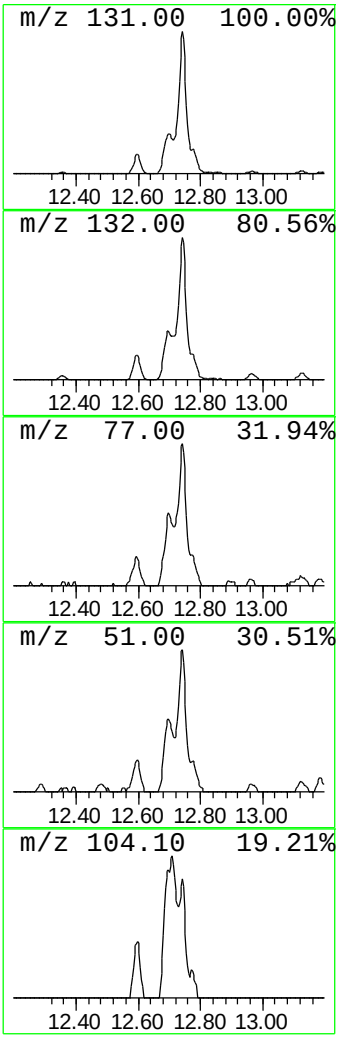
****

Peak Number 5 2-Propenal, 3-phenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	5.75 ug/L	65356	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	93
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
4	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	87
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

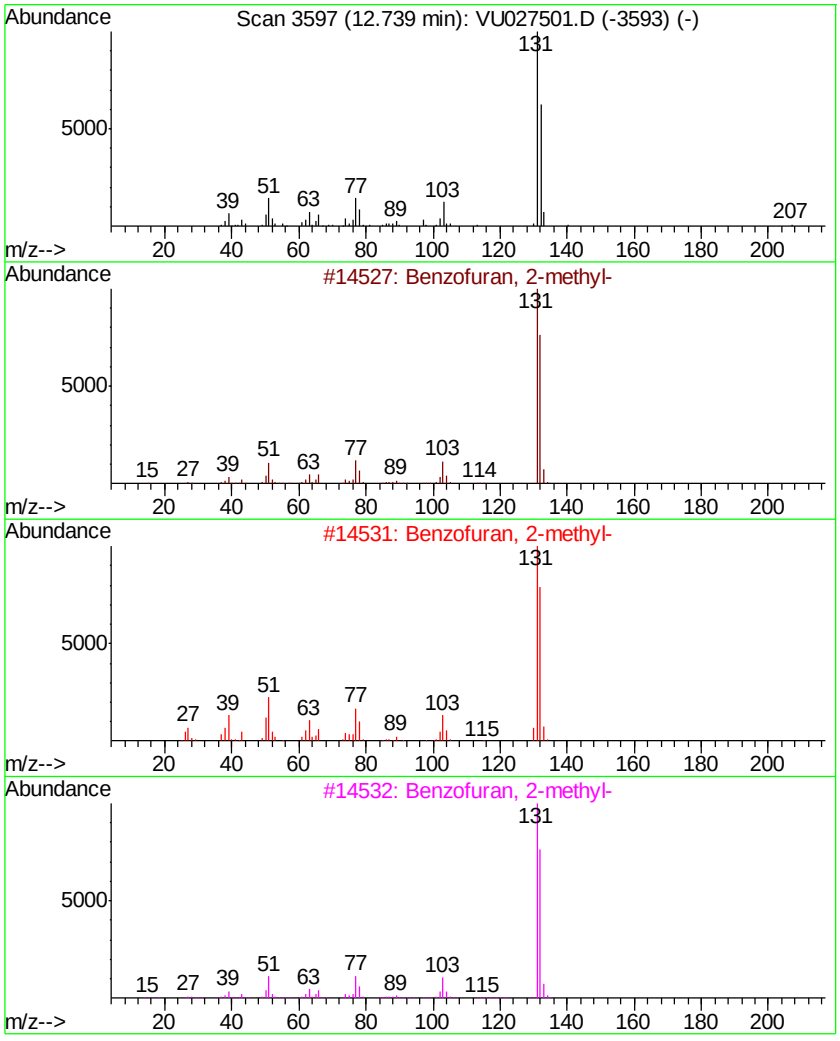
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

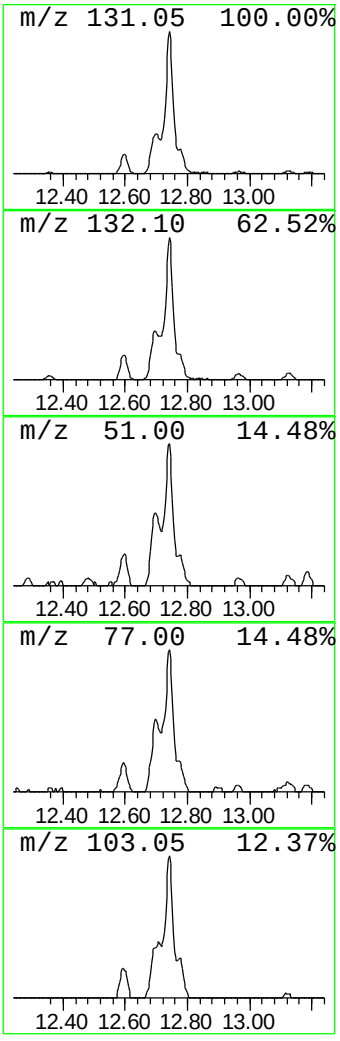
****

Peak Number 6 Benzofuran, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	13.59 ug/L	154382	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	86
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	86





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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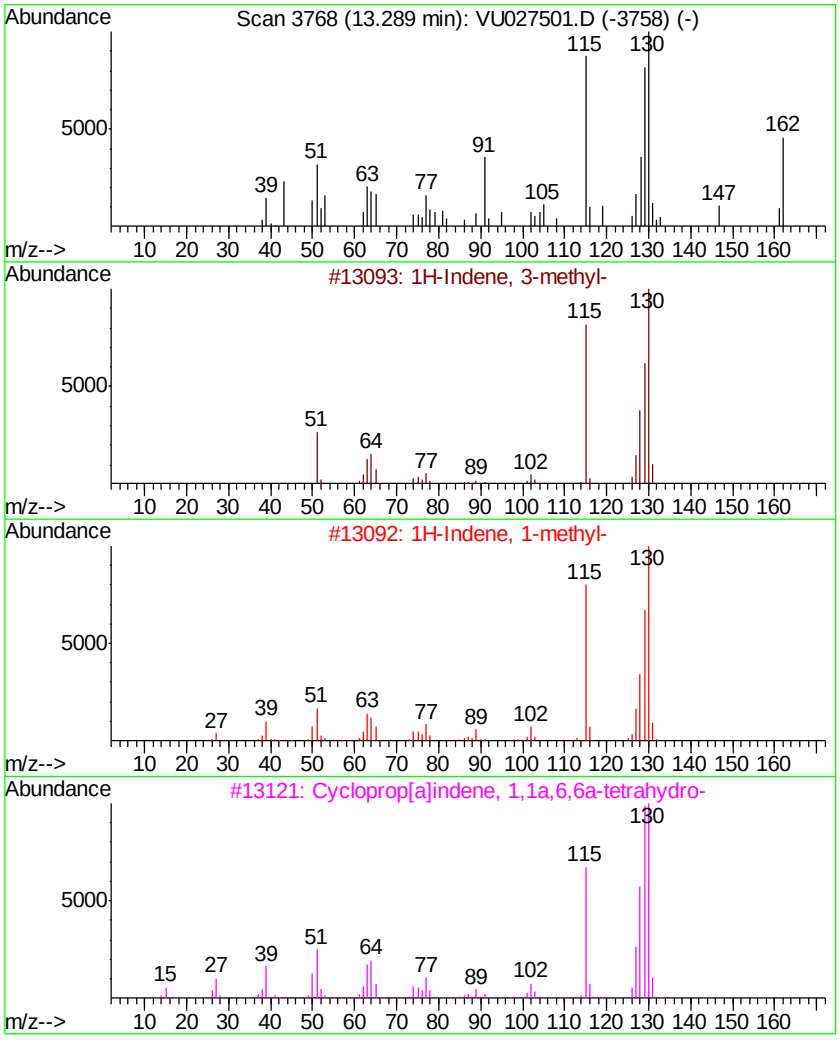
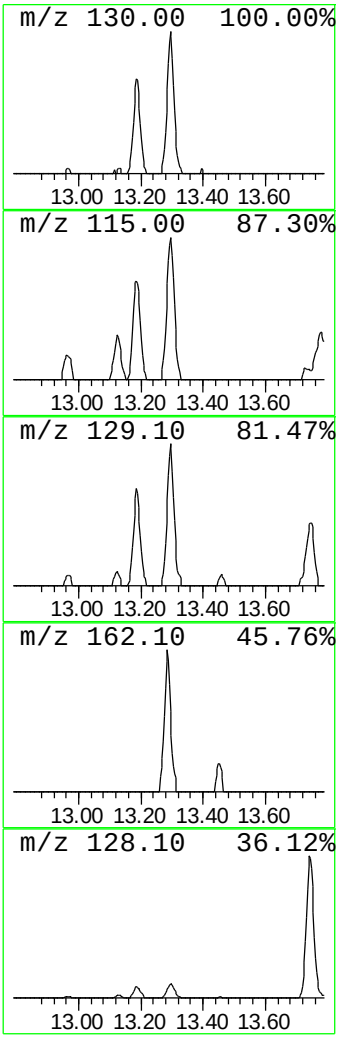
Peak Number 7 1H-Indene, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	3.27 ug/L	37130	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 3-methyl-	130	C10H10	000767-60-2	95
2	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
3	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	95
4	Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	93
5	2-Methylindene	130	C10H10	002177-47-1	93





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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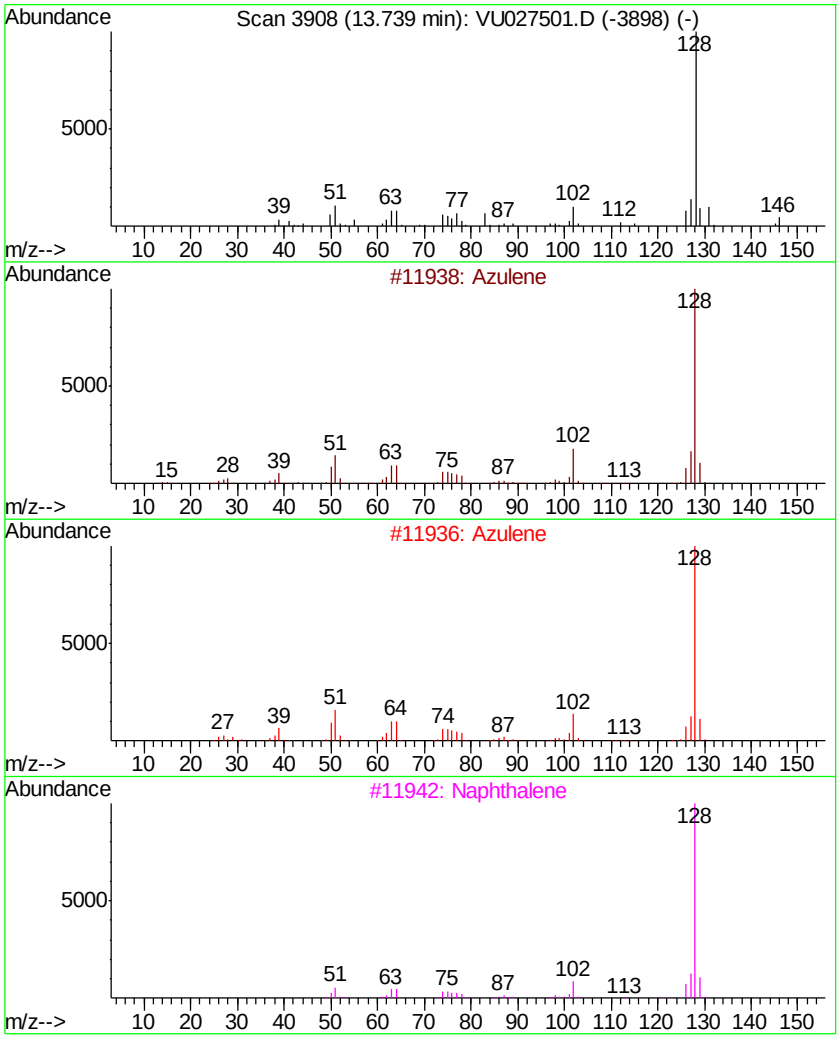
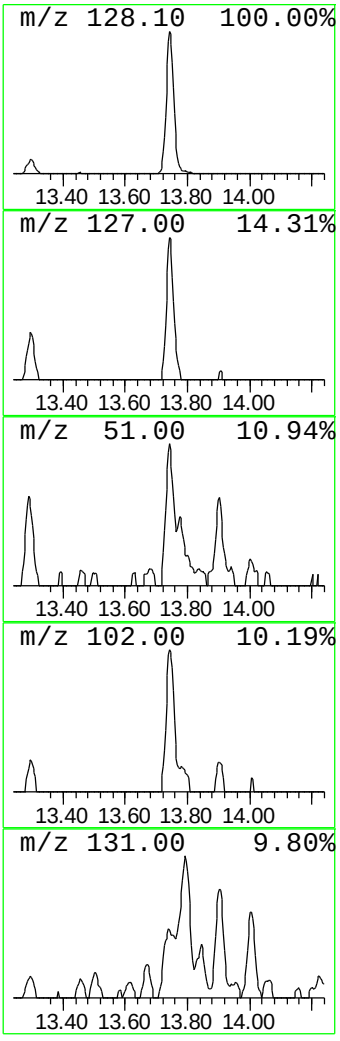
Peak Number 8 Azulene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.74	5.43 ug/L	61727	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Azulene	128	C10H8	000275-51-4	93
2	Azulene	128	C10H8	000275-51-4	93
3	Naphthalene	128	C10H8	000091-20-3	93
4	1H-Indene, 1-methylene-	128	C10H8	002471-84-3	93
5	Naphthalene	128	C10H8	000091-20-3	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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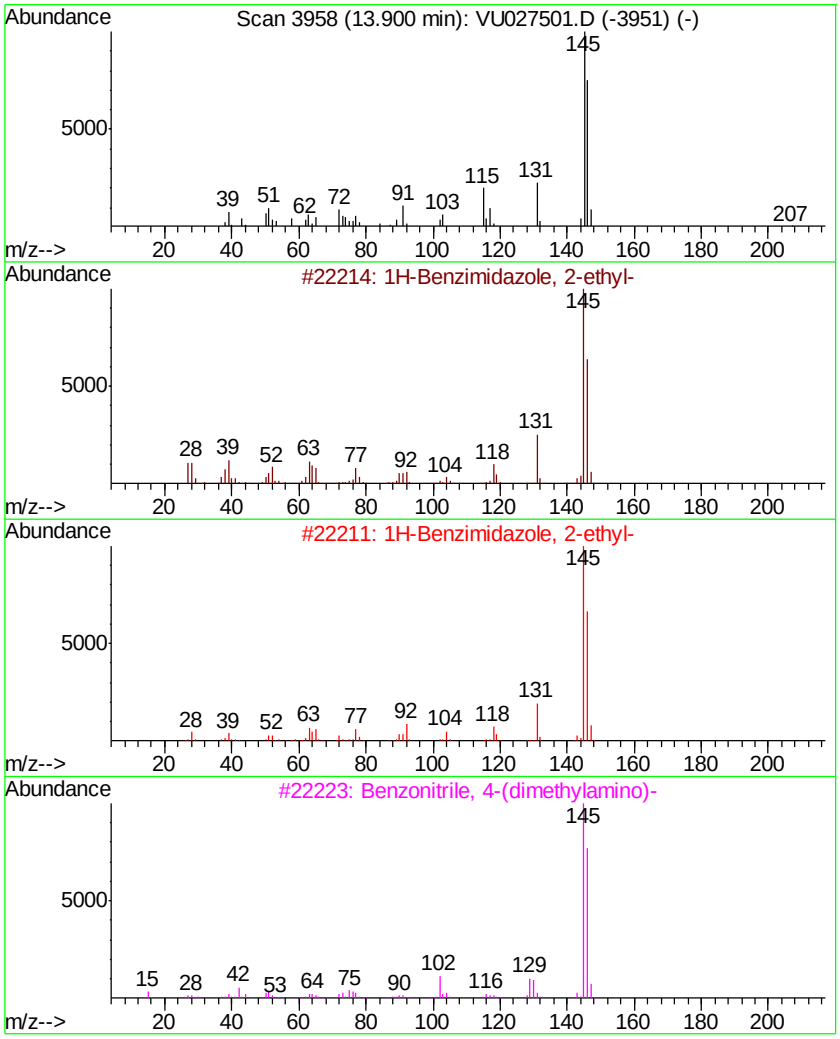
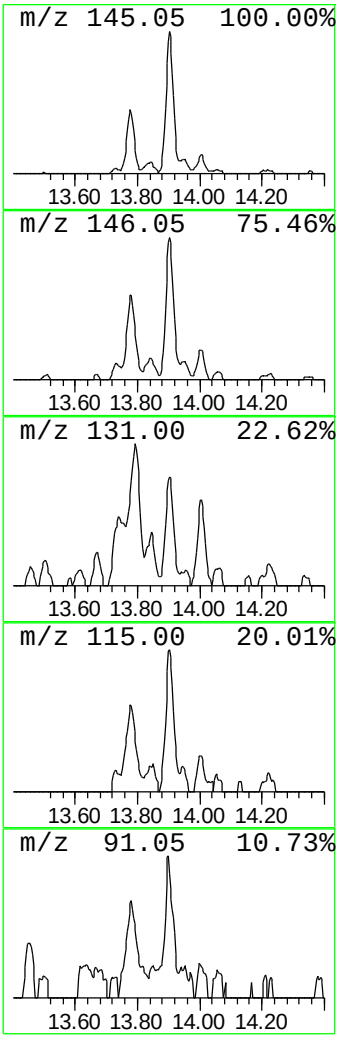
Peak Number 9 1H-Benzimidazole, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	3.73 ug/L	42376	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	64
2	1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	64
3	Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	53
4	1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	50
5	Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100818\
Data File : VU027501.D
Acq On : 08 Oct 2018 15:38
Operator : MD/SY
Sample : J5298-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : E62T6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM100518WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Indane	11.76	3.8	ug/L	43343	3	11.48	567924	50.0
Indene	11.98	3.4	ug/L	38102	3	11.48	567924	50.0
Benzofuran, 7-met...	12.59	2.6	ug/L	29422	3	11.48	567924	50.0
2-Propenal, 3-phe...	12.70	5.8	ug/L	65356	3	11.48	567924	50.0
Benzofuran, 2-met...	12.74	13.6	ug/L	154382	3	11.48	567924	50.0
1H-Indene, 3-methyl-	13.29	3.3	ug/L	37130	3	11.48	567924	50.0
Azulene	13.74	5.4	ug/L	61727	3	11.48	567924	50.0
1H-Benzimidazole,...	13.90	3.7	ug/L	42376	3	11.48	567924	50.0